Connecting via Winsock to STN

```
Welcome to STN International! Enter x:x
```

LOGINID:ssspta1712mxf

specific topic.

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TERMINAL (ENTER 1, 2, 3, OR ?):2

```
Welcome to STN International
                 Web Page URLs for STN Seminar Schedule - N. America
NEWS
                 "Ask CAS" for self-help around the clock
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                 New pricing for the Save Answers for SciFinder Wizard within
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                 KOREAPAT now available on STN
NEWS
         OCT 28
        NOV 30
                 PHAR reloaded with additional data
NEWS
      5
                 LISA now available on STN
        DEC 01
NEWS
        DEC 09
                 12 databases to be removed from STN on December 31, 2004
NEWS
      7
        DEC 15
                 MEDLINE update schedule for December 2004
NEWS 8
        DEC 17
                 ELCOM reloaded; updating to resume; current-awareness
NEWS
                 alerts (SDIs) affected
                 COMPUAB reloaded; updating to resume; current-awareness
      10 DEC 17
NEWS
                 alerts (SDIs) affected
                 SOLIDSTATE reloaded; updating to resume; current-awareness
NEWS
      11 DEC 17
                 alerts (SDIs) affected
                 CERAB reloaded; updating to resume; current-awareness
NEWS
      12 DEC 17
                 alerts (SDIs) affected
                 THREE NEW FIELDS ADDED TO IFIPAT/IFIUDB/IFICDB
      13 DEC 17
NEWS
                 EPFULL: New patent full text database to be available on STN
NEWS
      14 DEC 30
                 CAPLUS - PATENT COVERAGE EXPANDED
NEWS
      15 DEC 30
                 No connect-hour charges in EPFULL during January and
NEWS
      16 JAN 03
                 February 2005
                 CA/CAPLUS - Expanded patent coverage to include the Russian
NEWS
      17 JAN 26
                 Agency for Patents and Trademarks (ROSPATENT)
                 STN Patent Forums to be held in March 2005
NEWS
      18 FEB 10
              JANUARY 10 CURRENT WINDOWS VERSION IS V7.01a, CURRENT
NEWS EXPRESS
              MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP),
              AND CURRENT DISCOVER FILE IS DATED 10 JANUARY 2005
              STN Operating Hours Plus Help Desk Availability
NEWS HOURS
              General Internet Information
NEWS INTER
NEWS LOGIN
              Welcome Banner and News Items
              Direct Dial and Telecommunication Network Access to STN
NEWS PHONE
              CAS World Wide Web Site (general information)
NEWS WWW
```

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FILE 'HOME' ENTERED AT 16:10:04 ON 15 FEB 2005

=> file registry
COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 0.21 0.21

FULL ESTIMATED COST

FILE 'REGISTRY' ENTERED AT 16:10:10 ON 15 FEB 2005 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2005 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 14 FEB 2005 HIGHEST RN 831169-46-1 DICTIONARY FILE UPDATES: 14 FEB 2005 HIGHEST RN 831169-46-1

TSCA INFORMATION NOW CURRENT THROUGH MAY 21, 2004

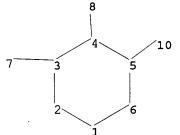
Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at: http://www.cas.org/ONLINE/DBSS/registryss.html

=>

Uploading C:\Program Files\Stnexp\Queries\10615694 la.str



chain nodes:
7 8 10
ring nodes:
1 2 3 4 5
chain bonds:
3-7 4-8 5-10
ring bonds:

1-2 1-6 2-3 3-4 4-5 5-6

exact/norm bonds :

1-2 1-6 2-3 3-4 3-7 4-5 4-8 5-6 5-10

G1:CH3,Et,H

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 10:CLASS

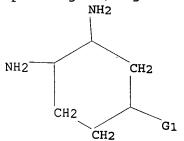
L1 STRUCTURE UPLOADED

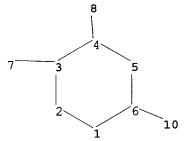
=> d l1 L1 HAS NO ANSWERS L1 STR

G1 Me,Et,H

Structure attributes must be viewed using STN Express query preparation.

Uploading C:\Program Files\Stnexp\Queries\10615694 lb.str





chain nodes :
7 8 10
ring nodes :
1 2 3 4 5 6
chain bonds :
3-7 4-8 6-10
ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6

exact/norm bonds :

1-2 1-6 2-3 3-4 3-7 4-5 4-8 5-6 6-10

G1:CH3,Et,H

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 10:CLASS

L2 STRUCTURE UPLOADED

=> d 12

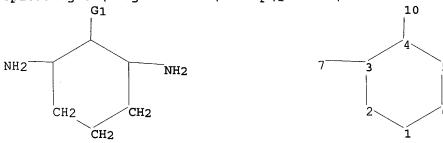
L2 HAS NO ANSWERS

L2 STR

G1 Me,Et,H

Structure attributes must be viewed using STN Express query preparation.

=>
Uploading C:\Program Files\Stnexp\Queries\10615694 1c.str



chain nodes :
7 8 10
ring nodes :
1 2 3 4 5 6
chain bonds :
3-7 4-10 5-8
ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6

exact/norm bonds :

1-2 1-6 2-3 3-4 3-7 4-5 4-10 5-6 5-8

G1:CH3,Et,H

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 10:CLASS

L3 STRUCTURE UPLOADED

=> d 13

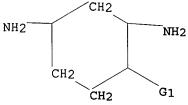
L3 HAS NO ANSWERS

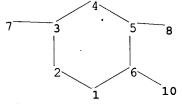
L3 STR

G1 Me,Et,H

Structure attributes must be viewed using STN Express query preparation.

=>
Uploading C:\Program Files\Stnexp\Queries\10615694 1d.str





chain nodes :
7 8 10
ring nodes :
1 2 3 4 5 6
chain bonds :
3-7 5-8 6-10
ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6

exact/norm bonds :

1-2 1-6, 2-3 3-4 3-7 4-5 5-6 5-8 6-10

G1:CH3,Et,H

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 10:CLASS

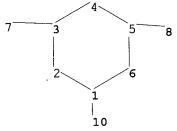
L4 STRUCTURE UPLOADED

=> d 14 L4 HAS NO ANSWERS L4 STR

G1 Me,Et,H

Structure attributes must be viewed using STN Express query preparation.

=>
Uploading C:\Program Files\Stnexp\Queries\10615694 1e.str



chain nodes :
7 8 10
ring nodes :
1 2 3 4 5
chain bonds :
1-10 3-7 5-8
ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6

exact/norm bonds :

1-2 1-6 1-10 2-3 3-4 3-7 4-5 5-6 5-8

G1:CH3,Et,H

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 10:CLASS

L5 STRUCTURE UPLOADED

=> d 15 L5 HAS NO ANSWERS

L5 STR

G1 Me, Et, H

Structure attributes must be viewed using STN Express query preparation.

Uploading C:\Program Files\Stnexp\Queries\10615694 1f.str

chain nodes:
7 8 10
ring nodes:
1 2 3 4 5 6
chain bonds:
1-10 3-7 6-8
ring bonds:

1-2 1-6 2-3 3-4 4-5 5-6

exact/norm bonds :

1-2 1-6 1-10 2-3 3-4 3-7 4-5 5-6 6-8

G1:CH3,Et,H

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 10:CLASS

L6 STRUCTURE UPLOADED

=> d 16 L6 HAS NO ANSWERS

L6 STR

G1 Me, Et, H

Structure attributes must be viewed using STN Express query preparation.

Uploading C:\Program Files\Stnexp\Queries\10615694 2a.str

chain nodes : 14 15 16 17 19 ring nodes : 2 3 4 5 7 8 9 10 11 12 13 6 chain bonds : 4-14 5-16 7-19 10-19 12-17 13-15 ring bonds : 4-5 5-6 6-7 8-9 8-13 9-10 10-11 11-12 12-13 2-3 2-7 3-4 exact/norm bonds : 2-3 2-7 3-4 4-5 4-14 5-6 5-16 6-7 8-9 8-13 9-10 10-11 11-12 12-13 12-17 13-15 exact bonds : 7-19 10-19

G1:CH3,Et,H

Match level:
2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:CLASS 15:CLASS 16:CLASS 17:CLASS 19:CLASS

L7 STRUCTURE UPLOADED

=> d 17 L7 HAS NO ANSWERS L7 STR

Structure attributes must be viewed using STN Express query preparation.

=>
Uploading C:\Program Files\Stnexp\Queries\10615694 2b.str

chain nodes : 14 15 16 17 19 ring nodes : 2 3 4 5 6 789 10 11 chain bonds : 4-14 5-16 7-19 10-19 12-17 13-15 ring bonds : 4-5 5-6 6-7 8-9 8-13 9-10 10-11 11-12 12-13 2-3 2-7 3-4 exact/norm bonds : $2-3 \quad 2-7 \quad 3-4 \quad 4-5 \quad 4-14 \quad 5-6 \quad 5-16 \quad 6-7 \quad 7-19 \quad 8-9 \quad 8-13 \quad 9-10 \quad 10-11 \quad 10-19$ 11-12 12-13 12-17 13-15

G1:CH3,Et,H

Match level:
2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:CLASS 15:CLASS 16:CLASS 17:CLASS 19:CLASS

L8 STRUCTURE UPLOADED

=> d 18 L8 HAS NO ANSWERS L8 STR

Structure attributes must be viewed using STN Express query preparation.

Uploading C:\Program Files\Stnexp\Queries\10615694 2c.str

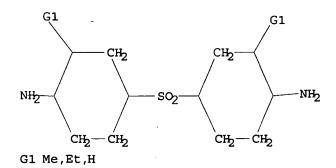
chain nodes : 14 15 16 17 ring nodes : 12 2 3 4 5 7 8 9 10 11 chain bonds : 4-14 5-16 7-19 10-19 12-17 13-15 ring bonds : 4-5 5-6 6-7 8-9 8-13 9-10 10-11 11-12 12-13 2-3 2-7 3-4 exact/norm bonds : 2-3 2-7 3-4 4-5 4-14 5-6 5-16 6-7 8-9 8-13 9-10 10-11 11-12 12-13 12-17 13-15 exact bonds : 7-19 10-19

G1:CH3,Et,H

Match level:
2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom
12:Atom 13:Atom 14:CLASS 15:CLASS 16:CLASS 17:CLASS 19:CLASS

L9 STRUCTURE UPLOADED

=> d 19 L9 HAS NO ANSWERS L9 STR



Structure attributes must be viewed using STN Express query preparation.

Uploading C:\Program Files\Stnexp\Queries\10615694 3a.str 24 NH2 CH2 CH2 $/19_{18}^{23}$ /CHCH2/ /5⁻6 /CHCH2 /1112 CH2 / NH144 _/716 NH2 / CH2 1710 /1315 3 2 CH2H2 9-8 CH2H2

chain nodes :
14 15 16 17 24

ring nodes :
2 3 4 5 6 7 8 9 10 11 12 13 18 19 20 21 22 23

chain bonds :
4-14 7-16 10-17 13-15 16-19 17-23 21-24

ring bonds :
2-3 2-7 3-4 4-5 5-6 6-7 8-9 8-13 9-10 10-11 11-12 12-13 18-19 18-23

19-20 20-21 21-22 22-23

exact/norm bonds :
2-3 2-7 3-4 4-5 4-14 5-6 6-7 8-9 8-13 9-10 10-11 11-12 12-13 13-15

18-19 18-23 19-20 20-21 21-22 21-24 22-23

exact bonds :
7-16 10-17 16-19 17-23

G1:CH3,Et,H

Match level:
2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom
12:Atom 13:Atom 14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS 19:Atom
20:Atom 21:Atom 22:Atom 23:Atom 24:CLASS

L10 STRUCTURE UPLOADED

L10 HAS NO ANSWERS L10 STR

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

Structure attributes must be viewed using STN Express query preparation.

Uploading C:\Program Files\Stnexp\Queries\10615694 3b.str

NH2 NH2 32 $20^{21}22$ $28^{29}30$ ĆH2 CH2 ĆH2 CH2 $/5^{-6}$ $/19_{18}^{23}$ $/27_{26}^{23}$ 1/CHCH2/ /CHCH2 CH2 /1112 CH_2 NH_1^{144} 7^{-16} NH2 / CH2 CH₂ 25 1710 /1315 CH2H2 9_8 CH2H2

chain nodes :

14 15 16 17 24 25 32

ring nodes :

2 3 4 5 6 7 8 9 10 11 12 13 18 19 20 21 22 23 26 27 28 29 30

31

chain bonds :

4-14 7-16 10-17 13-15 16-19 17-31 21-24 23-25 25-27 29-32

ring bonds :

2-3 2-7 3-4 4-5 5-6 6-7 8-9 8-13 9-10 10-11 11-12 12-13 18-19 18-23

19-20 20-21 21-22 22-23 26-27 26-31 27-28 28-29 29-30 30-31

exact/norm bonds :

2-3 2-7 3-4 4-5 4-14 5-6 6-7 8-9 8-13 9-10 10-11 11-12 12-13 13-15

18-19 18-23 19-20 20-21 21-22 21-24 22-23 26-27 26-31 27-28 28-29 29-30

29-32 30-31

exact bonds :

7-16 10-17 16-19 17-31 23-25 25-27

G1:CH3,Et,H

Match level:

2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom

12:Atom 13:Atom 14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS 19:Atom

20:Atom 21:Atom 22:Atom 23:Atom 24:CLASS 25:CLASS 26:CLASS 27:Atom 28:Atom

29:Atom 30:Atom 31:Atom 32:CLASS

STR

L11 STRUCTURE UPLOADED

=> d l11

L11 HAS NO ANSWERS

L11

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

Structure attributes must be viewed using STN Express query preparation.

```
=>
Uploading C:\Program Files\Stnexp\Queries\10615694 3c.str
          NH2
                                                        32
                 NH2
                        NH2
                              /CHCH2 /5<sup>-6</sup> /19/8<sup>23</sup> /27/31 /35/39

CH2 / NH\frac{144}{3} /7 16 25 33 CH2H2 3 2
                                                 20^{21}22 28^{29}30 36^{37}38
        CH2 CH2 CH2 CH2 CH2 CH2
                                                                         /1112
                         CH<sub>2</sub>
NH2 / CH2
                      CH<sub>2</sub>
                                                                       1710 /1315
  CH2H2
chain nodes :
14 15 16 17 24
                    25
                        32
                            33 40
ring nodes :
                             11 12 13 18 19 20 21 22 23
                                                                      27 28 29 30
           6789
                                                                 26
2 3 4 5
                        10
31 34 35 36 37 38
                        39
chain bonds :
4-14 7-16 10-17 13-15 16-19 17-39 21-24 23-25 25-27 29-32 31-33
                                                                             33-35
37-40
ring bonds :
2-3 2-7 3-4 4-5 5-6 6-7 8-9 8-13 9-10 10-11 11-12 12-13 18-19 18-23
19-20 20-21 21-22 22-23 26-27 26-31 27-28 28-29 29-30 30-31 34-35 34-39
35-36 36-37 37-38 38-39
exact/norm bonds :
2-3 2-7 3-4 4-5 4-14 5-6 6-7 8-9 8-13 9-10 10-11 11-12 12-13 13-15
18-19 18-23 19-20 20-21 21-22 21-24 22-23 26-27 26-31 27-28 28-29 29-30
29-32 30-31 34-35 34-39 35-36 36-37 37-38 37-40 38-39
exact bonds :
7-16 10-17 16-19 17-39 23-25 25-27 31-33 33-35
G1:CH3,Et,H
Match level :
2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom
12:Atom 13:Atom 14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS 19:Atom
20:Atom 21:Atom 22:Atom 23:Atom 24:CLASS 25:CLASS 26:CLASS 27:Atom 28:Atom 29:Atom 30:Atom 31:Atom 32:CLASS 33:CLASS 34:CLASS 35:Atom 36:Atom 37:Atom
38:Atom 39:Atom 40:CLASS
L12
        STRUCTURE UPLOADED
=> d 112
L12 HAS NO ANSWERS
                STR
*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***
```

Uploading C:\Program Files\Stnexp\Queries\10615694 3d.str

Structure attributes must be viewed using STN Express query preparation.

```
chain nodes :
14 15 16 17 24 25 32 33 40 41 48
ring nodes :
2 3 4 5 6 7 8 9 10 11 12 13 18 19 20 21 22 23 26 27 28 29 30
31 34 35 36 37 38 39 42 43 44 45 46 47
chain bonds :
4-14 7-16 10-17 13-15 16-19 17-47 21-24 23-25 25-27 29-32 31-33 33-35
37-40 39-41 41-43 45-48
ring bonds :
2-3 2-7 3-4 4-5 5-6 6-7 8-9 8-13 9-10 10-11 11-12 12-13 18-19 18-23
19-20 20-21 21-22 22-23 26-27 26-31 27-28 28-29 29-30 30-31 34-35 34-39
35-36 36-37 37-38 38-39 42-43 42-47 43-44 44-45 45-46 46-47
exact/norm bonds :
2-3 2-7 3-4 4-5 4-14 5-6 6-7 8-9 8-13 9-10 10-11 11-12 12-13 13-15
18-19 18-23 19-20 20-21 21-22 21-24 22-23 26-27 26-31 27-28 28-29 29-30 29-32 30-31 34-35 34-39 35-36 36-37 37-38 37-40 38-39 42-43 42-47 43-44 44-45 45-46 45-48 46-47
exact bonds :
7-16 10-17 16-19 17-47 23-25 25-27 31-33 33-35 39-41 41-43
```

G1:CH3,Et,H

Match level :

2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS 19:Atom 20:Atom 21:Atom 22:Atom 23:Atom 24:CLASS 25:CLASS 26:CLASS 27:Atom 28:Atom 29:Atom 30:Atom 31:Atom 32:CLASS 33:CLASS 34:CLASS 35:Atom 36:Atom 37:Atom 38:Atom 39:Atom 40:CLASS 41:CLASS 42:CLASS 43:Atom 44:Atom 45:Atom 46:Atom 47:Atom 48:CLASS

L13 STRUCTURE UPLOADED

Uploading C:\Program Files\Stnexp\Queries\10615694 4a.str

chain nodes :

14 15 16 17 24

ring nodes :

2 3 4 5 6 7 8 9 10 11 12 13 18 19 20 21 22 23

chain bonds :

4-14 7-16 10-17 13-15 16-19 17-23 21-24

ring bonds :

2-3 2-7 3-4 4-5 5-6 6-7 8-9 8-13 9-10 10-11 11-12 12-13 18-19 18-23

19-20 20-21 21-22 22-23

exact/norm bonds :

2-3 2-7 3-4 4-5 4-14 5-6 6-7 8-9 8-13 9-10 10-11 11-12 12-13 13-15

21-24

exact bonds :

7-16 10-17 16-19 17-23

normalized bonds :

18-19 18-23 19-20 20-21 21-22 22-23

G1:CH3,Et,H

Match level :

12:Atom 13:Atom 14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS 19:Atom

20:Atom 21:Atom 22:Atom 23:Atom 24:CLASS

L14 STRUCTURE UPLOADED

STR

=> d l14

L14 HAS NO ANSWERS

L14

$$\begin{array}{c} \text{CH}_2 - \text{CH}_2 \\ \end{array}$$

G1 Me, Et, H

Structure attributes must be viewed using STN Express query preparation.

```
Uploading C:\Program Files\Stnexp\Queries\10615694 4b.str
           NH2
                   NH2
                                                              22 28 30
          CH CH
                  CH CH
  /CHCH2<sup>/</sup>
                                                   /5<sup>-6</sup>
                                        /CHCH2
                                                                                          /1112
NH2 / CH2
                                      CH2 / NH141
                                                       /716
                                                                  25
                CH<sub>2</sub>
                                                                                        17L0 /1315
                                                    3-2
   CH2H2
                                         CH2H2
                                                                                           9 8
```

```
chain nodes :
14 15 16 17 24
ring nodes :
2 3 4 5
          6
             7
                  9
                     10
                         11
                            12
                               13
                                    18 19 20 21 22
                                                    23
                                                         26
                                                             27
                                                                28 29 30
31
chain bonds :
4-14 7-16 10-17 13-15 16-19 17-31
                                   21-24
                                          23-25 25-27
                                                      29-32
ring bonds :
2-3 2-7 3-4 4-5 5-6 6-7 8-9
                              8-13 9-10 10-11 11-12 12-13 18-19 18-23
19-20 20-21 21-22 22-23 26-27 26-31 27-28 28-29 29-30 30-31
exact/norm bonds :
2-3 2-7 3-4:4-5
                 4-14 5-6 6-7 8-9 8-13 9-10 10-11 11-12 12-13 13-15
21-24 29-32
exact bonds :
7-16 10-17 16-19 17-31 23-25 25-27
normalized bonds :
18-19 18-23 19-20 20-21 21-22 22-23 26-27 26-31 27-28 28-29 29-30 30-31
```

G1:CH3,Et,H

```
Match level:
2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom
12:Atom 13:Atom 14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS 19:Atom
20:Atom 21:Atom 22:Atom 23:Atom 24:CLASS 25:CLASS 26:CLASS 27:Atom 28:Atom
29:Atom 30:Atom 31:Atom 32:CLASS
```

L15 STRUCTURE UPLOADED

=> d 115 L15 HAS NO ANSWERS L15 STR

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

Structure attributes must be viewed using STN Express query preparation.

=> Uploading C:\Program Files\Stnexp\Queries\10615694 4c.str NH2 NH2 NH2 CH CH CH CH CH /CHCH2/ CH₂CH /1112 /CHCH2 CH CH2 / NH141 /716 CH2H2 3 2 NH2 / CH2 17L0 /1315 CH2H2 9 8

chain nodes : 14 15 16 17 24 25 32 33 ring nodes : 2 3 4 5 6 7 8 9 10 11 12 13 18 19 20 21 22 23 26 27 28 29 30 31 34 35 36 37 38 39 chain bonds : 4-14 7-16 10-17 13-15 16-19 17-39 21-24 23-25 25-27 29-32 31-33 33-35 37-40 ring bonds : 2-3 2-7 3-4 4-5 5-6 6-7 8-9 8-13 9-10 10-11 11-12 12-13 18-19 18-23 19-20 20-21 21-22 22-23 26-27 26-31 27-28 28-29 29-30 30-31 34-35 34-39 35-36 36-37 37-38 38-39 exact/norm bonds : 2-3 2-7 3-4 4-5 4-14 5-6 6-7 8-9 8-13 9-10 10-11 11-12 12-13 13-15 21-24 29-32 37-40 exact bonds : 7-16 10-17 16-19 17-39 23-25 25-27 31-33 33-35 normalized bonds : $18 - 19 \quad 18 - 23 \quad 19 - 20 \quad 20 - 21 \quad 21 - 22 \quad 22 - 23 \quad 26 - 27 \quad 26 - 31 \quad 27 - 28 \quad 28 - 29 \quad 29 - 30 \quad 30 - 31$ 34-35 34-39 35-36 36-37 37-38 38-39

G1:CH3,Et,H

Match level :
2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom
12:Atom 13:Atom 14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS 19:Atom
20:Atom 21:Atom 22:Atom 23:Atom 24:CLASS 25:CLASS 26:CLASS 27:Atom 28:Atom
29:Atom 30:Atom 31:Atom 32:CLASS 33:CLASS 34:CLASS 35:Atom 36:Atom 37:Atom
38:Atom 39:Atom 40:CLASS

L16 STRUCTURE UPLOADED

=> d 116 L16 HAS NO ANSWERS L16 STR

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

Structure attributes must be viewed using STN Express query preparation.

```
Uploading C:\Program Files\Stnexp\Queries\10615694 4d.str
                                           24
                                                32
        NH2
              NH2
                                                       40
                    NH2 NH2
                                    CH CH CH CH CH CHCH CH
                    CH
 /CHCH2/
                           /CHCH2
                         CH/
           CH<sub>2</sub>CH
                          CH2 / NH141 /716
NH2 / CH2
                                                              17L0 /1:15
                  CH2
                      CH2
  CH2H2
                                                                9 8
                             CH2H2
```

```
chain nodes :
14 15 16 17 24 25
                      32
                          33
                              40
                                  41
                                      48
ring nodes :
                                                  21 22 23 26 27 28 29 30
2 3 4 5
          6 7 8 9
                       10
                              12
                                  13
                                      18
                                         19 20
                          11
          36 37 38
31 34 35
                      39
                          42
                              43
                                  44
                                      45
                                         46 47
chain bonds :
4-14 7-16 10-17 13-15 16-19 17-47 21-24 23-25 25-27 29-32 31-33 33-35
37-40 39-41 41-43 45-48
ring bonds :
2-3 \quad 2-7 \quad 3-4 \quad 4-5 \quad 5-6 \quad 6-7 \quad 8-9 \quad 8-13 \quad 9-10 \quad 10-11 \quad 11-12 \quad 12-13 \quad 18-19 \quad 18-23
19-20 20-21 21-22 22-23 26-27 26-31 27-28 28-29 29-30 30-31 34-35 34-39
35-36 36-37 37-38 38-39 42-43 42-47 43-44 44-45 45-46 46-47
exact/norm bonds :
2-3 2-7 3-4 4-5 4-14 5-6 6-7 8-9 8-13 9-10 10-11 11-12 12-13 13-15
21-24 29-32 37-40 45-48
exact bonds :
7-16 10-17 16-19 17-47 23-25 25-27 31-33 33-35 39-41 41-43
normalized bonds :
18-19 18-23 19-20 20-21 21-22 22-23 26-27 26-31 27-28 28-29 29-30 30-31
34-35 34-39 35-36 36-37 37-38 38-39 42-43 42-47 43-44 44-45 45-46 46-47
```

G1:CH3,Et,H

```
Match level:
2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom
12:Atom 13:Atom 14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS 19:Atom
20:Atom 21:Atom 22:Atom 23:Atom 24:CLASS 25:CLASS 26:CLASS 27:Atom 28:Atom
29:Atom 30:Atom 31:Atom 32:CLASS 33:CLASS 34:CLASS 35:Atom 36:Atom 37:Atom
38:Atom 39:Atom 40:CLASS 41:CLASS 42:CLASS 43:Atom 44:Atom 45:Atom 46:Atom
47:Atom 48:CLASS
```

=> d 117

L17 HAS NO ANSWERS

L17

STR

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

Structure attributes must be viewed using STN Express query preparation.

=> s sam (11 or 12 or 13 or 14 or 15 or 16 or 17 or 18 or 19 or 110 or 111 or 112

or 113 or 114 or 115 or 116 or 117)

SAMPLE SEARCH INITIATED 16:19:12 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 61276 TO ITERATE

1.6% PROCESSED 1000 ITERATIONS 3 ANSWERS

INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **INCOMPLETE**

BATCH **INCOMPLETE**

PROJECTED ITERATIONS: EXCEEDS 1000000

PROJECTED ANSWERS: EXCEEDS 2863

L18 3 SEA SSS SAM (L1 OR L2 OR L3 OR L4 OR L5 OR L6 OR L7 OR L8 OR L9

=> s sam l1

SAMPLE SEARCH INITIATED 16:19:29 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 15382 TO ITERATE

6.5% PROCESSED 1000 ITERATIONS 2 ANSWERS

INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS: 300213 TO 315067

PROJECTED ANSWERS: 283 TO 947

L19 2 SEA SSS SAM L1

=> s sam 12

SAMPLE SEARCH INITIATED 16:19:34 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 15382 TO ITERATE

6.5% PROCESSED 1000 ITERATIONS 2 ANSWERS

INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

1 ANSWERS

PROJECTED ITERATIONS: 300213 TO 315067

PROJECTED ANSWERS: 283 TO 947

L20 2 SEA SSS SAM L2

=> s sam 13

SAMPLE SEARCH INITIATED 16:19:39 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED - 21368 TO ITERATE

4.7% PROCESSED 1000 ITERATIONS

INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **INCOMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS: 418613 TO 436107

PROJECTED ANSWERS: 150 TO 704

L21 1 SEA SSS SAM L3

=> s sam 14 SAMPLE SEARCH INITIATED 16:19:43 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED - 21368 TO ITERATE

4.7% PROCESSED 1000 ITERATIONS 1 ANSWERS

INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **INCOMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS: 418613 TO 436107

150 TO PROJECTED ANSWERS: 704

1 SEA SSS SAM L4 L22

=> s sam 15 SAMPLE SEARCH INITIATED 16:19:46 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED - 21368 TO ITERATE

4.7% PROCESSED 1000 ITERATIONS 1 ANSWERS

INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **INCOMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS: 418613 TO 436107

PROJECTED ANSWERS: 150 TO 704

1 SEA SSS SAM L5 L23

=> s sam 16 SAMPLE SEARCH INITIATED 16:19:50 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED - 28012 TO ITERATE

3.6% PROCESSED 1000 ITERATIONS 0 ANSWERS

INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **INCOMPLETE**

BATCH **COMPLETE**

550233 TO 570247 PROJECTED ITERATIONS: 0 TO PROJECTED ANSWERS:

0 SEA SSS SAM L6 L24

=> s sam 17 SAMPLE SEARCH INITIATED 16:19:57 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED - 6473 TO ITERATE

15.4% PROCESSED 1000 ITERATIONS 13 ANSWERS INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS: 124637 TO 134283

PROJECTED ANSWERS: 1132 TO 2232

L25 13 SEA SSS SAM L7

=> s sam 18 SAMPLE SEARCH INITIATED 16:20:01 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED - 4857 TO ITERATE

0 ANSWERS 20.6% PROCESSED 1000 ITERATIONS

INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS: 92961 TO 101319

PROJECTED ANSWERS: 0 TO

0 SEA SSS SAM L8

=> s sam 19 SAMPLE SEARCH INITIATED 16:20:04 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED - 723 TO ITERATE

100.0% PROCESSED 723 ITERATIONS 2 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

12847 TO 16073 PROJECTED ITERATIONS:

PROJECTED ANSWERS: 2 TO 124

2 SEA SSS SAM L9

=> s sam 110 SAMPLE SEARCH INITIATED 16:20:08 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED - 27 TO ITERATE

0 ANSWERS 100.0% PROCESSED 27 ITERATIONS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**

229 TO 851 PROJECTED ITERATIONS: PROJECTED ANSWERS: O TO

0 SEA SSS SAM L10

=> s sam l11 SAMPLE SEARCH INITIATED 16:20:11 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED - 17 TO ITERATE

100.0% PROCESSED 17 ITERATIONS 0 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS: 93 TO 587 PROJECTED ANSWERS: 0 TO

L29 0 SEA SSS SAM L11

=> s sam 112

SAMPLE SEARCH INITIATED 16:20:14 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 4 TO ITERATE

100.0% PROCESSED 4 ITERATIONS 0 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

PROJECTED ITERATIONS: 4 TO 200

PROJECTED ANSWERS: 0 TO 0

L30 0 SEA SSS SAM L12

=> s sam 113

SAMPLE SEARCH INITIATED 16:20:17 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 8 TO ITERATE

100.0% PROCESSED 8 ITERATIONS 0 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 8 TO 329

PROJECTED ANSWERS: 0 TO 0

L31 0 SEA SSS SAM L13

=> s sam 114

SAMPLE SEARCH INITIATED 16:20:22 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 20 TO ITERATE

100.0% PROCESSED 20 ITERATIONS 0 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**

PROJECTED ITERATIONS: 132 TO 668
PROJECTED ANSWERS: 0 TO 0

L32 0 SEA SSS SAM L14

=> s sam l15

SAMPLE SEARCH INITIATED 16:20:26 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 12 TO ITERATE

100.0% PROCESSED 12 ITERATIONS 0 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**

PROJECTED ITERATIONS: 33 TO 447
PROJECTED ANSWERS: 0 TO 0

L33 0 SEA SSS SAM L15

=> s sam 116

SAMPLE SEARCH INITIATED 16:20:28 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 3 TO ITERATE

100.0% PROCESSED 3 ITERATIONS 0 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS: 3 TO 163

PROJECTED ANSWERS: O TO

O SEA SSS SAM L16

=> s sam 117

SAMPLE SEARCH INITIATED 16:20:32 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED - 4 TO ITERATE

100.0% PROCESSED 4 ITERATIONS 0 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS: 4 TO 200 PROJECTED ANSWERS: 0 TO

L35 0 SEA SSS SAM L17

=> s full l1

FULL SEARCH INITIATED 16:20:34 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 310971 TO ITERATE

100.0% PROCESSED 310971 ITERATIONS 390 ANSWERS

SEARCH TIME: 00.00.02

L36 390 SEA SSS FUL L1

=> s full 12

FULL SEARCH INITIATED 16:20:41 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 310971 TO ITERATE

100.0% PROCESSED 310971 ITERATIONS 421 ANSWERS

SEARCH TIME: 00.00.02

L37 421 SEA SSS FUL L2

=> s full 13

FULL SEARCH INITIATED 16:20:47 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 429126 TO ITERATE

93.2% PROCESSED 400000 ITERATIONS 131 ANSWERS

INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.03

FULL FILE PROJECTIONS: ONLINE **INCOMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS: 429126 TO 429126

131 TO PROJECTED ANSWERS: 175

131 SEA SSS FUL L3 L38

=> s full 14

FULL SEARCH INITIATED 16:20:53 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 429126 TO ITERATE

93.2% PROCESSED 400000 ITERATIONS 183 ANSWERS

INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.03

FULL FILE PROJECTIONS: ONLINE **INCOMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS: 429126 TO 429126 238

PROJECTED ANSWERS: 183 TO

183 SEA SSS FUL L4 L39

=> s full 15

FULL SEARCH INITIATED 16:21:00 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 429126 TO ITERATE

93.2% PROCESSED 400000 ITERATIONS 218 ANSWERS

INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.03

FULL FILE PROJECTIONS: ONLINE **INCOMPLETE**
BATCH **COMPLETE**

PROJECTED ITERATIONS: 429126 TO 429126

218 TO PROJECTED ANSWERS: 278

218 SEA SSS FUL L5 L40

=> s full 16

FULL SEARCH INITIATED 16:21:20 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 563587 TO ITERATE

71.0% PROCESSED 400000 ITERATIONS 220 ANSWERS

INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.04

FULL FILE PROJECTIONS: ONLINE **INCOMPLETE**
BATCH **COMPLETE**

PROJECTED ITERATIONS: 563587 TO 563587

PROJECTED ANSWERS: 257 TO 361

L41 220 SEA SSS FUL L6

=> s full 17

FULL SEARCH INITIATED 16:21:28 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 130033 TO ITERATE

100.0% PROCESSED 130033 ITERATIONS 1517 ANSWERS

SEARCH TIME: 00.00.01

L42 1517 SEA SSS FUL L7

=> s full 18

FULL SEARCH INITIATED 16:21:33 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 96964 TO ITERATE

100.0% PROCESSED 96964 ITERATIONS 33 ANSWERS

SEARCH TIME: 00.00.01

33 SEA SSS FUL L8 L43

=> s full 19

FULL SEARCH INITIATED 16:21:37 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 14211 TO ITERATE

100.0% PROCESSED 14211 ITERATIONS 5 ANSWERS

SEARCH TIME: 00.00.01

5 SEA SSS FUL L9 L44

=> s full 110

FULL SEARCH INITIATED 16:21:44 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 359 TO ITERATE

100.0% PROCESSED 359 ITERATIONS 0 ANSWERS

SEARCH TIME: 00.00.01

L45 0 SEA SSS FUL L10

=> s full 111

FULL SEARCH INITIATED 16:21:50 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 292 TO ITERATE

100.0% PROCESSED 292 ITERATIONS 0 ANSWERS

SEARCH TIME: 00.00.01

L46 0 SEA SSS FUL L11

=> s full 112

FULL SEARCH INITIATED 16:21:53 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 86 TO ITERATE

100.0% PROCESSED 86 ITERATIONS 0 ANSWERS

SEARCH TIME: 00.00.01

L47 0 SEA SSS FUL L12

=> s full 113

FULL SEARCH INITIATED 16:21:58 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 182 TO ITERATE

100.0% PROCESSED 182 ITERATIONS 0 ANSWERS

SEARCH TIME: 00.00.01

L48 0 SEA SSS FUL L13

=> s full 114

FULL SEARCH INITIATED 16:22:01 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 294 TO ITERATE

100.0% PROCESSED 294 ITERATIONS 0 ANSWERS

SEARCH TIME: 00.00.01

L49 0 SEA SSS FUL L14

=> s full 115

FULL SEARCH INITIATED 16:22:06 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 253 TO ITERATE

100.0% PROCESSED 253 ITERATIONS 0 ANSWERS

SEARCH TIME: 00.00.01

L50 0 SEA SSS FUL L15

=> s full 116

FULL SEARCH INITIATED 16:22:08 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 140 TO ITERATE

100.0% PROCESSED 140 ITERATIONS 0 ANSWERS

SEARCH TIME: 00.00.01

L51 0 SEA SSS FUL L16

=> s full 117

FULL SEARCH INITIATED 16:22:13 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 174 TO ITERATE

100.0% PROCESSED 174 ITERATIONS 0 ANSWERS

SEARCH TIME: 00.00.01

L52 0 SEA SSS FUL L17

=> file caplus

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION

FULL ESTIMATED COST 2743.90 2744.11

FILE 'CAPLUS' ENTERED AT 16:22:24 ON 15 FEB 2005 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2005 AMERICAN CHEMICAL SOCIETY (ACS)

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FILE COVERS 1907 - 15 Feb 2005 VOL 142 ISS 8 FILE LAST UPDATED: 14 Feb 2005 (20050214/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s (136 or 137 or 138 or 139 or 140 or 141 or 142 or 143 or 144 or 145 or 146 or 147 or 148 or 149 or 150 or 151 or 152)

2037 L36

2035 L37

137 L38

154 L39

253 L40

182 L41

2673 L42

12 L43

3 L44

0 L45

0 L46

0 L47 0 L48

0 L49

0 L50

0 L51

0 T₁52

L53 5038 (L36 OR L37 OR L38 OR L39 OR L40 OR L41 OR L42 OR L43 OR L44 OR L45 OR L46 OR L47 OR L48 OR L49 OR L50 OR L51 OR L52)

=> s 153 and (ketone or ketimine or diketimine or polyketimine)

137872 KETONE

112328 KETONES

200999 KETONE

(KETONE OR KETONES)

```
10615694 2/15/05
```

```
2092 KETIMINE
          1303 KETIMINES .
          2611 KETIMINE
                 (KETIMINE OR KETIMINES)
           172 DIKETIMINE
            61 DIKETIMINES
           204 DIKETIMINE
                 (DIKETIMINE OR DIKETIMINES)
            36 POLYKETIMINE
            19 POLYKETIMINES
            42 POLYKETIMINE
                 (POLYKETIMINE OR POLYKETIMINES)
           202 L53 AND (KETONE OR KETIMINE OR DIKETIMINE OR POLYKETIMINE)
L54
=> 154 and (epox? or polyepox? or diepox? or diglycidyl? or polyglycidyl? or
glycidylether or glycidylether?)
L54 IS NOT A RECOGNIZED COMMAND
The previous command name entered was not recognized by the system.
For a list of commands available to you in the current file, enter
"HELP COMMANDS" at an arrow prompt (=>).
=> s (154 and (epox? or polyepox? or diepox? or diglycidyl? or polyglycidyl? or
glycidylether or glycidylether?))
        294154 EPOX?
          4365 POLYEPOX?
          5861 DIEPOX?
         16833 DIGLYCIDYL?
          2189 POLYGLYCIDYL?
            32 GLYCIDYLETHER
             2 GLYCIDYLETHERS
            34 GLYCIDYLETHER
                  (GLYCIDYLETHER OR GLYCIDYLETHERS)
            36 GLYCIDYLETHER?
            27 (L54 AND (EPOX? OR POLYEPOX? OR DIEPOX? OR DIGLYCIDYL? OR POLYGL
L55
               YCIDYL? OR GLYCIDYLETHER OR GLYCIDYLETHER?))
=> s 155 and (ketimine or diketimine or triketimine or polyketimine)
          2092 KETIMINE
          1303 KETIMINES
          2611 KETIMINE
                  (KETIMINE OR KETIMINES)
           172 DIKETIMINE
            61 DIKETIMINES
           204 DIKETIMINE
                  (DIKETIMINE OR DIKETIMINES)
             3 TRIKETIMINE
             1 TRIKETIMINES
             4 TRIKETIMINE
                  (TRIKETIMINE OR TRIKETIMINES)
            36 POLYKETIMINE
            19 POLYKETIMINES
            42 POLYKETIMINE
                  (POLYKETIMINE OR POLYKETIMINES)
             6 L55 AND (KETIMINE OR DIKETIMINE OR TRIKETIMINE OR POLYKETIMINE)
L56
=> d 156 abs ibib hitstr 1
L56 ANSWER 1 OF 6 CAPLUS COPYRIGHT 2005 ACS on STN
     The composition comprises a polyepoxide compound and a ketimine
     derived from the condensation reaction of an amine having \geq 2
     primary amino groups directed bonded to a cyclohexane ring and an aliphatic
     ketone. Thus, a composition comprising ketimine prepared from
     1,2-diaminocyclohexane and 4-methyl-2-pentanone 6.7, xylene 3 and
```

bisphenol A diglycidyl ether 19 parts showed initial viscosity

165 mPa-s, and 14 days required for twice higher the initial viscosity and

40 days required for three times higher the initial viscosity.

ACCESSION NUMBER: 2005:34641 CAPLUS

DOCUMENT NUMBER: 142:115129

TITLE: Curable epoxy resin compositions for

adhesives, sealants and coatings with longer pot life INVENTOR(S): Chiba, Yasuo; Morimoto, Hiroya; Yokomakura, Masahiro

PATENT ASSIGNEE(S): Japan

SOURCE: U.S. Pat. Appl. Publ., 9 pp.

CODEN: USXXCO

DOCUMENT TYPE:

LANGUAGE:

Patent English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.				KINI)	DATE		P	APPL:	ICAT:	ION I	NO.		D	ATE				
							-			-						_			
	US	2005	0100	22		A1		2005	0113	τ	JS 20	003-	6156	94		20	0030	709	
	ΕP	1496	075			A1		2005	0112	E	EP 20	004-	1594	9		2	040	707	
		R:	•	•		•		•		GB,			•		•		-	_	
			ΙE,	SI,	LT,	LV,	FΙ,	RO,	MK,	CY,	ΑL,	TR,	ВG,	CZ,	EE,	HU,	PL,	SK,	HR
	JP	2005	0297	97		A2		2005	0203	Ċ	JP 20	004-	2034	20		2	040	709	
PRIORITY APPLN. INFO.:					Ţ	JS 2	003-	6156	94	7	A 2	0030	709						
TM COA 02 7DD 1 2 Disminosyral chowspo					2020	~+ i ~:	n nr	04110	+0 10	ith									

IT 694-83-7DP, 1,2-Diaminocyclohexane, reaction products with

ketones

RL: IMF (Industrial manufacture); RCT (Reactant); PREP (Preparation); RACT (Reactant or reagent)

(curing agent; curable **epoxy** resin compns. for adhesives, sealants and coatings with longer pot life)

RN 694-83-7 CAPLUS

CN 1,2-Cyclohexanediamine (6CI, 7CI, 8CI, 9CI) (CA INDEX NAME)

=> d l56 abs ibib hitstr 2

L56 ANSWER 2 OF 6 CAPLUS COPYRIGHT 2005 ACS on STN

AB The title amines, giving cured compns. forming clear films, are prepared from compds. bearing ≥2 NH2 groups, alkylene carbonates, aliphatic carbonyl compds., and compds. bearing ≥2 OH-reactive groups in specified ratios. Adding 1.0 mol ethylene carbonate over 1 h to 1.0 mol 1,6-hexanediamine in PhMe at 40-50°, heating at 50° until the amine number was <158, adding 2.0 mol MEK and 0.5 g MeC6H4SO3H, distilling H2O, adding 0.5 mol polypropylene glycol-based diepoxy resin (epoxy content 3125 mmol/kg), and heating at 120° until the epoxy content was <50 mmol/kg gave a blocked amine. Use of 76 g this product with 50 g aqueous epoxy resin (Beckopox EP 116) composition (epoxy content 5076 mmol/kg) in the formulation of a clear coating film is exemplified.

ACCESSION NUMBER: 1999:736300 CAPLUS

DOCUMENT NUMBER: 131:337884

TITLE: Blocked amines as curing agent for aqueous

one-component epoxy resin compositions

INVENTOR(S): Feola, Roland; Muller, Friedrich; Gmoser, Johann

PATENT ASSIGNEE(S): Vianova Resins AG, Austria; Surface Specialties

Austria GM

SOURCE: Eur. Pat. Appl., 11 pp.

CODEN: EPXXDW

DOCUMENT TYPE:

Patent

LANGUAGE:

י. ז

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PA?	TENT	NO.			KIN	D DATE		A	PL	ICAT:	ION I	NO.		D	ATE	
														_		
EP	9571	21			A2	1999	1117	E	1:	999-:	1089	40		1	9990	505
EP	9571	21			A3	2002	0522									
EP	9571	21			B1	2004	0204									
	R:	ΑT,	BE,	CH,	DE,	DK, ES,	FR,	GB, G	R,	IT,	LI,	LU,	NL,	SE,	MC,	PT,
		ΙE,	SI,	LT,	LV,	FI, RO										
AT	9800	802			Α	2000	1015	ΑT	1:	998-1	802			1	9980	512
AT	4077	48			В	2001	0525									
AT	2589	44			E	2004	0215	A	1:	999-:	1089	40		1	9990	505
US	6207	733			· B1	2001	0327	US	1	999-3	3079	74		1	9990	510
PRIORIT	Y APP	LN.	INFO	. :				A7	1:	998-	802			A 1	9980	512
								EI	1	999-	1089	40		A 1	9990	505

IT 694-83-7D, 1,2-Cyclohexanediamine, reaction products with alkylene carbonates, **ketones** and **epoxy** resins

1761-71-3D, Bis(4-aminocyclohexyl) methane, reaction products with

alkylene carbonates, ketones and epoxy resins

RL: MOA (Modifier or additive use); USES (Uses)

(blocked amines as curing agent for aqueous one-component epoxy resin compns.)

RN 694-83-7 CAPLUS

CN 1,2-Cyclohexanediamine (6CI, 7CI, 8CI, 9CI) (CA INDEX NAME)

RN 1761-71-3 CAPLUS

CN Cyclohexanamine, 4,4'-methylenebis- (9CI) (CA INDEX NAME)

=> d 156 abs ibib hitstr 3

L56 ANSWER 3 OF 6 CAPLUS COPYRIGHT 2005 ACS on STN

AB Title agents are obtained by reaction of (A) ketone resins with (B) amines H2N[R1(NH2)pNR2(R3NR4)m]R5 [R1 = (aromatic or heterocyclic ring-containing) C1-32 linear or branched alkylene, (alkyl-substituted) cycloalkylene; R2, R4 = H, C1-4 linear or branched alkyl; R3 = (OH-substituted) alkylene; R5 = H, (aromatic or heterocyclic ring-terminated) alkyl, cycloalkyl, cycloalkylalkylene; p, m = 0-1; n = 0-18] with a mol ratio of NH2 (amine) to CO (ketone resin) 1/3 to 3/1. Thus, reacting 136 g m-xylenediamine with 150 g an acetophenone-based

ketone resin (CO equivalent 150) gave 150 g a compound, 32 parts of which were blended with 100 parts EP 4100 (an **epoxy** resin) to show tack-free time 5 h at 23°, pot life 12 h, and good oil- and weather- resistance when applied on a soft steel plate.

ACCESSION NUMBER: 1996:672175 CAPLUS

DOCUMENT NUMBER: 125:277661

TITLE: Crosslinking agents with long pot life and short

crosslinking time for epoxy resins

INVENTOR(S): Akimoto, Koji; Hayano, Satoshi; Kusano, Shoji;

Tsukada, Yasutada

PATENT ASSIGNEE(S): Asahi Denka Kogyo KK, Japan

SOURCE: Jpn. Kokai Tokkyo Koho, 9 pp.

CODEN: JKXXAF

DOCUMENT TYPE: Patent LANGUAGE: Japanese

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 08217858	A2	19960827	JP 1995-29780	19950217
JP 3526946	B2	20040517		
D.T 3.D.D.T.17 T.1700			TD 1005 00500	

PRIORITY APPLN. INFO.: JP 1995-29780 19950217

IT 1761-71-3DP, Bis(4-aminocyclohexyl)methane, reaction products with ketone resins

RL: IMF (Industrial manufacture); RCT (Reactant); PREP (Preparation); RACT (Reactant or reagent)

(crosslinking agents; **ketimine** crosslinking agents with long pot life for **epoxy** resins)

RN 1761-71-3 CAPLUS

CN Cyclohexanamine, 4,4'-methylenebis- (9CI) (CA INDEX NAME)

=> d 156 abs ibib hitstr 4

L56 ANSWER 4 OF 6 CAPLUS COPYRIGHT 2005 ACS on STN AB The title resins, dispersible in H2O on protonation and useful in electrodip coatings, contain (A) reaction products of epoxy resins [number-average mol. weight (Mn) 300-6000, 1.5-3.0 epoxy groups/mol.] with diketimines from diprimary amines (and optionally a primary monoamine ketimine, a secondary amine optionally containing a tertiary amino group, and/or a ketimine of a primary/tertiary diamine) and (B) reaction products of epoxy resins (Mn 800-6000) with secondary amines (optionally bearing tertiary amino groups), ketimines of primary monoamines, and/or ketimines of primary-tertiary diamines. A mixture of 640 g reaction product of bisphenol A diglycidyl ether (I) 214.3, bisphenol A 48, 387:580:97 1,6-hexanediamine-dimer acid-linseed-oil fatty acid reaction product iso-BuCOMe ketimine 210.9, MeNHCH2CH2OH 3.22, and diethanolamine (II) 18 g; 233.3 g reaction product of I 121.3 and II 26.3 g; 322 g reaction product of trimethylolpropane 1340, urea 3600, Bu2NH 7740, 1,6-hexanediamine 1740, and 4,4'-methylenebis(2methylcyclohexylamine) 3570 g in aqueous AcOH gave a pigmented coating (deposited at 320 V and baked 20 min at 170°) with impact deepening (ASTM D 2794) 18.08 N-m, surface appearance (1 best, 6 worst) 1, and

undercutting in climate cycling (VDA 621415) 1.5 mm.

ACCESSION NUMBER: 1989:214828 CAPLUS

DOCUMENT NUMBER: 110:214828

TITLE: Resins containing basic nitrogen for waterborne

coatings

INVENTOR(S): Perner, Thomas; Osterloh, Rolf; Schupp, Eberhard;

Schwerzel, Thomas; Ahlers, Klaas

PATENT ASSIGNEE(S): BASF Lacke und Farben A.-G., Fed. Rep. Ger.

SOURCE: Eur. Pat. Appl., 11 pp.

CODEN: EPXXDW

DOCUMENT TYPE: Patent LANGUAGE: German

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PAC	TENT NO.		KIND	DATE	APPLICATION NO.		DATE
EP	296494		A2	19881228	EP 1988-109640		19880616
EP	296494		A3	19900516			
EP	296494		B1	19930908			
	R: AT,	BE, CH	I, DE, 1	ES, FR, GB,	IT, LI, NL, SE		
DE	3720955		A1	19890105	DE 1987-3720955		19870625
AΤ	94183		E	19930915	AT 1988-109640		19880616
ES	2058184		Т3	19941101	ES 1988-109640		19880616
US	4981884		Α	19910101	US 1988-210502		19880623
BR	8803108		Α	19890124	BR 1988-3108		19880624
JP	01022955		A2	19890125	JP 1988-155108		19880624
ZA	8804527		Α	19900228	ZA 1988-4527		19880624
PRIORIT	Y APPLN.	INFO.:			DE 1987-3720955	Α	19870625
					EP 1988-109640	Α	19880616

IT 120658-95-9

RL: MOA (Modifier or additive use); USES (Uses) (crosslinking agents, for electrophoretic coatings)

RN 120658-95-9 CAPLUS

CN Urea, polymer with N-butyl-1-butanamine, 2-ethyl-2-(hydroxymethyl)-1,3-propanediol, 1,6-hexanediamine and 4,4'-methylenebis[2-methylcyclohexanamine] (9CI) (CA INDEX NAME)

CM 1

CRN 6864-37-5 CMF C15 H30 N2

$$\begin{array}{c} \text{Me} \\ \\ \text{H}_2 \text{N} \\ \end{array} \begin{array}{c} \text{NH}_2 \\ \\ \text{Me} \end{array}$$

CM 2

CRN 124-09-4 CMF C6 H16 N2

 $H_2N-(CH_2)_6-NH_2$

CM 3

CRN 111-92-2 CMF C8 H19 N

n-Bu-NH-Bu-n

CM 4

CRN 77-99-6 CMF C6 H14 O3

$${\rm CH_2-OH} \ {\rm Ho-CH_2-C-Et} \ {\rm CH_2-OH} \ {\rm CH_2-O$$

CM 5

CRN 57-13-6 CMF C H4 N2 O

$$H_2N-C-NH_2$$

=> d 156 abs ibib hitstr 5

L56 ANSWER 5 OF 6 CAPLUS COPYRIGHT 2005 ACS on STN

The title liquid materials contain (meth)acryloyl compound and C2-24 (cyclo)aliphatic primary amines [or their adducts (mol. weight 350-1500) with epoxides, isocyanates, or unsatd. carbonyl compds.] blocked by C≤10 aldehydes or ketones. Thus, a polymer solution (from xylene 3640, glycidyl methacrylate 1560, styrene 2807, and Bu acrylate 1872 g) 4252, acrylic acid 320.4, Cr(III) 2-ethylhexanoate 2.7, and hydroquinone 4.57 g were mixed, aerated at 110° to acid value <2, and mixed with 150 g xylene to give a 53.1% solution with acid value 0.9 and viscosity 170 cPa-s. An equiequiv. mixture of this solution and blocked amine (from dipropylenetriamine 196.5, iso-BuCOMe 330, and Epikote-828 297.7 g) was coated to 40 μ (dry basis) on steel to give a coating with tack-free time 2 h, Persoz hardness 104 and 213 after 1 and 7 days, and good resistance to MEK and gasoline.

ACCESSION NUMBER:

1987:215612 CAPLUS

DOCUMENT NUMBER:

106:215612

TITLE:

Fast-drying discoloration-free coating materials

AKZO N. V., Neth.

PATENT ASSIGNEE(S): SOURCE:

Jpn. Kokai Tokkyo Koho, 11 pp.

CODEN: JKXXAF

DOCUMENT TYPE:

Patent

LANGUAGE:

Japanese

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

			APPLICATION NO.	
JP 61252273			JP 1986-68829	
JP 61252273 JP 05053186	B4 .	19930809		
			EP 1986-103844	19860320
EP 203296				
EP 203296	B2	19970903		
R: AT, BE, CH,				
AT 46712	E	19891015	AT 1986-103844	19860320
DK 8601385	Α	19860930	DK 1986-1385	19860325
DK 172448	B1	19980713		
NO 8601222	A	19860930	NO 1986-1222	19860325
NO 175864	В	19940912		
NO 175864	C	19941221		
BR 8601354	A	19861202	BR 1986-1354	
ES 553476	A1	19870601	ES 1986-553476	19860326
CA 1268889	A1	19900508	CA 1986-505213	19860326
CN 86101964	Α	19860924	CN 1986-101964	19860327
	В			
FI 8601346			FI 1986-1346	19860327
	В			
FI 80468		19900611		
AU 8655366	A1	19861002	AU 1986-55366	19860327
AU 584689	B2	19890601		
ZA 8602359	A	19861126		
US 4990577	A	19910205	US 1989-309449	19890210
PRIORITY APPLN. INFO.:			NL 1985-952	
			EP 1986-103844	A 19860320
			US 1986-844412	B1 19860326

IT **6864-37-5D**, 3,3'-Dimethyl-4,4'-diaminodicyclohexylmethane, reaction products with **ketones**

RL: USES (Uses)

(crosslinking agents for room temperature-curable acrylic coatings)

RN 6864-37-5 CAPLUS

CN Cyclohexanamine, 4,4'-methylenebis[2-methyl- (9CI) (CA INDEX NAME)

$$\begin{array}{c} \text{Me} \\ \\ \text{H}_2 \text{N} \\ \\ \text{Me} \end{array}$$

=> d 156 abs ibib hitstr 6

L56 ANSWER 6 OF 6 CAPLUS COPYRIGHT 2005 ACS on STN

The reaction of an epoxy resin with a diketimine of a compound containing 2 primary amino groups and, optionally, with a secondary amine (optionally containing a tertiary amino group), a ketimine of a primary monoamine, and/or a ketimine of primary-tertiary diamine gives a resin which contains basic N-containing groups. The resin are treated with an acid and used in the preparation of water-thinned compns. suitable for cathodic deposition in the preparation of coatings. The cured coatings have good solvent resistance and salt-spray resistance. Thus, H2N(CH2)6MH2 232, dimerized fatty acids 290, and iso-BuCOMe 621 parts were used to prepared a ketimine. The reaction of 210.9 parts of this ketimine and 22.5 parts diethanolamine with 207.9 part bisphenol A-epichlorohydrin copolymer (I) having epoxide equivalent 485 and

54.3 parts I having epoxide equivalent 190 gave a resin containing basic nitrogen-containing groups. This resin, a blocked isocyanate, AcOH, and dibutyltin dilaurate were used with a pigment paste in preparation of coating compns. which were coated on Zn phosphate-treated steel by cathodic deposition and cured 20 min at 180°.

ACCESSION NUMBER:

1985:150998 CAPLUS

DOCUMENT NUMBER:

102:150998

TITLE:

Synthetic resin containing basic nitrogen-containing

groups and its use

INVENTOR (S):

Schupp, Eberhard; Loch, Werner; Osterloh, Rolf;

Ahlers, Klaas

PATENT ASSIGNEE(S):

BASF Farben und Fasern A.-G., Fed. Rep. Ger.

SOURCE:

Ger. Offen., 17 pp. CODEN: GWXXBX

DOCUMENT TYPE:

Patent

LANGUAGE:

German

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.		DATE
				-	
DE 3325061	A1	19850124	DE 1983-3325061		19830712
EP 134983	A1	19850327	EP 1984-107958		19840706
EP 134983	B1	19870603			
R: AT, BE, CH,	DE, FR	, GB, IT, L	I, NL, SE		
AT 27608	E	19870615	AT 1984-107958		19840706
ZA 8405328	Α	19850327	ZA 1984-5328		19840710
BR 8403432	Α	19850625	BR 1984-3432		19840710
JP 60063223	A2	19850411	JP 1984-142507		19840711
JP 04049843	B4	19920812			
ES 534228	A1	19850416	ES 1984-534228	•	19840711
US 4557814	Α	19851210	US 1984-629762		19840711
CA 1232099	A1	19880126	CA 1984-458661		19840711
PRIORITY APPLN. INFO.:			DE 1983-3325061	A	19830712
			EP 1984-107958	Α	19840706

IT 6864-37-5D, reaction products with urea

RL: USES (Uses)

(hardeners, for **epoxy** resin-ketimine reaction

products in coatings)

RN 6864-37-5 CAPLUS

Cyclohexanamine, 4,4'-methylenebis[2-methyl- (9CI) (CA INDEX NAME) CN

$$\begin{array}{c|c} \text{Me} & \\ \text{H}_2\text{N} & \\ \text{Me} & \\ \end{array}$$

=> s 155 and ketone and blocked and (polyamine or diamine or triamine)

137872 KETONE

112328 KETONES

200999 KETONE

(KETONE OR KETONES)

137757 BLOCKED

1 BLOCKEDS

137757 BLOCKED

(BLOCKED OR BLOCKEDS)

31476 POLYAMINE

```
10615694 2/15/05
```

33319 POLYAMINES 45012 POLYAMINE

(POLYAMINE OR POLYAMINES)

41487 DIAMINE 25587 DIAMINES

57405 DIAMINE

(DIAMINE OR DIAMINES)

4026 TRIAMINE 852 TRIAMINES

4607 TRIAMINE

(TRIAMINE OR TRIAMINES)

3 L55 AND KETONE AND BLOCKED AND (POLYAMINE OR DIAMINE OR TRIAMINE L57

=> d 157 abs ibib hitstr 1

L57 ANSWER 1 OF 3 CAPLUS COPYRIGHT 2005 ACS on STN GI

$$\begin{array}{c|c}
0 & x \\
\hline
 & R1
\end{array}$$

Ι

AB Blocked polyurethanes that form reactive systems with organic compds. having ≥2 amine groups and, optionally, epoxides having >1 epoxide group are manufactured by reaction of ≥1 polyisocyanate with ≥1 polyol and blocking the excess NCO groups with activated ketones I [X = electron-withdrawing group, R1, R2 = H or C≥12 (un)saturated, or (cyclo)aliphatic groups and (substituted) aromatic or (substituted) araliph. groups, optionally containing ≤3 0, S, or N atoms, n = 0-5] in the optionally presence of a catalyst. These blocked polyurethanes have lower viscosity and form cured products with no low-mol.-weight compds. resulting from the deblocking during curing. A typical blocked polyurethane was manufactured by rapidly adding 74.17 g 2,4-tolylene diisocyanate to 852.58 g Acclaim 4200 at 60°, stirring at 60° until the NCO content is 1.93%, adding 1 g Zn 2-ethylhexanoate and 73.25 g Et cyclopentanone-2-carboxylate, and stirring at 50° until the NCO content was 0.1%.

ACCESSION NUMBER:

2004:525942 CAPLUS

DOCUMENT NUMBER:

141:89883

TITLE:

Reactive blocked polyurethanes

INVENTOR(S):

Simon, Joachim; Karlou-eyrisch, Kamelia; Guertler, Christoph; Mager, Michael; Schelhaas, Michael; Stingl,

Thomas

PATENT ASSIGNEE(S):

Bayer Ag, Germany

Ger. Offen., 14 pp. SOURCE:

CODEN: GWXXBX

DOCUMENT TYPE:

Patent

LANGUAGE:

German

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PATENT NO. APPLICATION NO. KIND DATE DATE

```
A1
                                20040701
                                            DE 2002-10260299
                                                                   20021220
    DE 10260299
    WO 2004058849
                         A1
                                20040715
                                            WO 2003-EP13833
                                                                   20031206
        W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH,
            CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD,
            GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC,
            LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NI, NO,
            NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ,
            TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW
        RW: BW, GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ,
            BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE,
            ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK,
            TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD,
                                20040729
                                            US 2003-738742
                                                                   20031217
    US 2004147704
                         A1
                                                                A 20021220
PRIORITY APPLN. INFO.:
                                            DE 2002-10260299
     713541-91-4P, Acclaim 4200-ethyl cyclopentanone-2-carboxylate-
     Laromin C260-2,4-tolylene diisocyanate copolymer 713541-92-5P,
     Acclaim 2200-ethyl cyclopentanone-2-carboxylate-Laromin C260-2,4-tolylene
     diisocyanate copolymer 713541-93-6P, Acclaim 1000-ethyl
     cyclopentanone-2-carboxylate-Laromin C260-2,4-tolylene diisocyanate
     copolymer 713541-94-7P, Acclaim 3201-ethyl cyclopentanone-2-
     carboxylate-Laromin C260-2,4-tolylene diisocyanate copolymer
     713541-95-8P, Ethyl cyclopentanone-2-carboxylate-ethylene
     oxide-Laromin C260-2,4-propylene oxide-tolylene diisocyanate copolymer
     713541-96-9P, Acclaim 8200-ethyl cyclopentanone-2-carboxylate-
     Laromin C260-2,4-tolylene diisocyanate copolymer 713541-97-0P,
     Acclaim 12200-ethyl cyclopentanone-2-carboxylate-Laromin C260-2,4-tolylene
     diisocyanate copolymer 713541-98-1P, Acclaim 2200-ethyl
     cyclohexanone-2-carboxylate-Laromin C260-2,4-tolylene diisocyanate
     copolymer 713541-99-2P, Acclaim 4200-ethyl cyclohexanone-2-
     carboxylate-Laromin C260-2,4-tolylene diisocyanate copolymer
     713542-00-8P, Ethyl cyclohexanone-2-carboxylate-Laromin
     C260-polypropylene glycol trimethylolpropane ether-TDI copolymer
     714248-43-8P, Ethylene oxide-propylene oxide copolymer glycerol
     ether-ethyl cyclopentanone-2-carboxylate-Laromin C260-2,4-tolylene
     diisocyanate copolymer
     RL: IMF (Industrial manufacture); PREP (Preparation)
        (cured sample; reactive blocked polyurethanes prepared with
        activated cyclic ketones as blocking agents)
     713541-91-4 CAPLUS
RN
     Cyclopentanecarboxylic acid, 2-oxo-, ethyl ester, polymer with Acclaim
CN
     4200, 2,4-diisocyanato-1-methylbenzene and 4,4'-methylenebis[2-
     methylcyclohexanamine] (9CI) (CA INDEX NAME)
     CM
          188571-36-0
     CRN
     CMF
          Unspecified
     CCI
          PMS, MAN
*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***
     CM
     CRN
          6864-37-5
          C15 H30 N2
     CMF
```

$$\begin{array}{c} \text{Me} \\ \text{H}_2\text{N} \\ \end{array} \begin{array}{c} \text{CH}_2 \\ \text{Me} \\ \end{array}$$

CM 3

CRN 611-10-9 CMF C8 H12 O3

CM 4

CRN 584-84-9 CMF C9 H6 N2 O2

RN 713541-92-5 CAPLUS

CN Cyclopentanecarboxylic acid, 2-oxo-, ethyl ester, polymer with Acclaim 2200, 2,4-diisocyanato-1-methylbenzene and 4,4'-methylenebis[2-methylcyclohexanamine] (9CI) (CA INDEX NAME)

CM 1

CRN 188571-34-8 CMF Unspecified CCI PMS, MAN

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

CM 2

$$\begin{array}{c} \text{Me} \\ \\ \text{H}_2 \text{N} \\ \end{array} \begin{array}{c} \text{CH}_2 \\ \\ \text{Me} \\ \end{array}$$

3 CM

611-10-9 CRN CMF C8 H12 O3

CM

CRN 584-84-9 CMF C9 H6 N2 O2

713541-93-6 CAPLUS RN

Cyclopentanecarboxylic acid, 2-oxo-, ethyl ester, polymer with Acclaim 1000, 2,4-diisocyanato-1-methylbenzene and 4,4'-methylenebis[2-methylcyclohexanamine] (9CI) (CA INDEX NAME) CN

CM

CRN 694464-11-4 Unspecified CMF CCI PMS, MAN

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

CM

$$\begin{array}{c} \text{Me} \\ \\ \text{H}_2 \text{N} \\ \end{array} \begin{array}{c} \text{CH}_2 \\ \\ \text{Me} \\ \end{array}$$

CM 3

CRN 611-10-9 CMF C8 H12 O3

CM 4

CRN 584-84-9 CMF C9 H6 N2 O2

RN 713541-94-7 CAPLUS

CN Cyclopentanecarboxylic acid, 2-oxo-, ethyl ester, polymer with Acclaim 3201, 2,4-diisocyanato-1-methylbenzene and 4,4'-methylenebis[2-methylcyclohexanamine] (9CI) (CA INDEX NAME)

CM 1

CRN 188571-35-9 CMF Unspecified CCI PMS, MAN

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

CM 2

$$\begin{array}{c} \text{Me} \\ \text{H}_2\text{N} \\ \end{array} \begin{array}{c} \text{CH}_2 \\ \text{Me} \\ \end{array}$$

CM 3

CRN 611-10-9 CMF C8 H12 O3

CM 4

CRN 584-84-9 CMF C9 H6 N2 O2

RN 713541-95-8 CAPLUS

CN Cyclopentanecarboxylic acid, 2-oxo-, ethyl ester, polymer with 2,4-diisocyanato-1-methylbenzene, 4,4'-methylenebis[2-methylcyclohexanamine] and oxirane (9CI) (CA INDEX NAME)

CM 1

CRN 6864-37-5 CMF C15 H30 N2

$$\begin{array}{c} \text{Me} \\ \\ \text{H}_2 \text{N} \\ \end{array} \begin{array}{c} \text{CH}_2 \\ \\ \text{Me} \\ \end{array}$$

CM 2

CRN 611-10-9

CMF C8 H12 O3

CM 3

CRN 584-84-9 C9 H6 N2 O2 CMF

. CM

CRN 75-21-8 CMF C2 H4 O



CN

RN713541-96-9 CAPLUS

Cyclopentanecarboxylic acid, 2-oxo-, ethyl ester, polymer with Acclaim 8200, 2,4-diisocyanato-1-methylbenzene and 4,4'-methylenebis[2-methylcyclohexanamine] (9CI) (CA INDEX NAME)

CM

CRN 188571-38-2 CMF Unspecified CCI PMS, MAN

STRUCTURE DIAGRAM IS NOT AVAILABLE ***

CM

6864-37-5 CRN C15 H30 N2 CMF

$$\begin{array}{c} \text{Me} \\ \\ \text{H}_2\text{N} \\ \end{array} \begin{array}{c} \text{CH}_2 \\ \\ \text{Me} \\ \end{array}$$

3 CM

CRN 611-10-9 CMF C8 H12 O3

CM 4 .

584-84-9 CRN C9 H6 N2 O2 CMF

RN713541-97-0 CAPLUS

Cyclopentanecarboxylic acid, 2-oxo-, ethyl ester, polymer with Acclaim 12200, 2,4-diisocyanato-1-methylbenzene and 4,4'-methylenebis[2-methylcyclohexanamine] (9CI) (CA INDEX NAME) CN

CM

CRN 278793-29-6 CMF Unspecified CCI PMS, MAN

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

CM 2

$$H_2N$$
 CH_2
 NH_2
 Me

CM 3

CRN 611-10-9 CMF C8 H12 O3

CM 4

CRN 584-84-9 CMF C9 H6 N2 O2

RN 713541-98-1 CAPLUS

CN Cyclohexanecarboxylic acid, 2-oxo-, ethyl ester, polymer with Acclaim 2200, 2,4-diisocyanato-1-methylbenzene and 4,4'-methylenebis[2-methylcyclohexanamine] (9CI) (CA INDEX NAME)

CM 1

CRN 188571-34-8 CMF Unspecified CCI PMS, MAN

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

CM 2

$$\begin{array}{c} \text{Me} \\ \text{H}_2\text{N} \\ \end{array} \begin{array}{c} \text{CH}_2 \\ \text{Me} \end{array}$$

CM 3

CRN 1655-07-8 CMF C9 H14 O3

CM 4

CRN 584-84-9 CMF C9 H6 N2 O2

RN 713541-99-2 CAPLUS

CN Cyclohexanecarboxylic acid, 2-oxo-, ethyl ester, polymer with Acclaim 4200, 2,4-diisocyanato-1-methylbenzene and 4,4'-methylenebis[2-methylcyclohexanamine] (9CI) (CA INDEX NAME)

CM 1

CRN 188571-36-0 CMF Unspecified CCI PMS, MAN

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

CM 2

$$\begin{array}{c} \text{Me} \\ \\ \text{H}_2 \text{N} \\ \\ \text{Me} \end{array}$$

CM 3

CRN 1655-07-8 CMF C9 H14 O3

CM 4

CRN 584-84-9 CMF C9 H6 N2 O2

RN 713542-00-8 CAPLUS

CN Cyclohexanecarboxylic acid, 2-oxo-, ethyl ester, polymer with 1,3-diisocyanatomethylbenzene, α-hydro-ω-hydroxypoly[oxy(methyl-1,2-ethanediyl)] ether with 2-ethyl-2-(hydroxymethyl)-1,3-propanediol (3:1), and 4,4'-methylenebis[2-methylcyclohexanamine] (9CI) (CA INDEX NAME)

CM 1

CRN 26471-62-5 CMF C9 H6 N2 O2 CCI IDS

CM 2

CRN 25723-16-4

CMF (C3 H6 O)n (C3 H6 O)n (C3 H6 O)n C6 H14 O3

CCI IDS, PMS

$$\begin{array}{c|c} CH_2 & \hline & O-(C_3H_6) & \hline & \\ HO & \hline & \\ CH_2 & \hline & O-(C_3H_6) & \hline \\ CH_2 &$$

CM 3

CRN 6864-37-5 CMF C15 H30 N2

$$\begin{array}{c} \text{Me} \\ \\ \text{H}_2 \text{N} \\ \end{array} \begin{array}{c} \text{CH}_2 \\ \\ \text{Me} \\ \end{array}$$

CM 4

CRN 1655-07-8 CMF C9 H14 O3

RN 714248-43-8 CAPLUS

CN Cyclopentanecarboxylic acid, 2-oxo-, ethyl ester, polymer with 2,4-diisocyanato-1-methylbenzene, 4,4'-methylenebis[2-methylcyclohexanamine] and methyloxirane polymer with oxirane ether with 1,2,3-propanetriol (3:1) (9CI) (CA INDEX NAME)

CM 1

$$\begin{array}{c} \text{Me} \\ \\ \text{H}_2 \text{N} \\ \\ \text{Me} \\ \end{array}$$

CM 2

CRN 611-10-9 CMF C8 H12 O3

CM 3

CRN 584-84-9 CMF C9 H6 N2 O2

CM 4

CRN 9082-00-2

CMF C3 H8 O3 . 3 (C3 H6 O . C2 H4 O) x

CM 5

CRN 56-81-5 CMF C3 H8 O3

$$\begin{array}{c} \text{OH} \\ | \\ \text{HO- CH}_2\text{- CH- CH}_2\text{- OH} \end{array}$$

CM 6

CRN 9003-11-6

CMF (C3 H6 O . C2 H4 O) x

CCI PMS

CM 7

CRN 75-56-9 C3 H6 O CMF



CM 8

CRN 75-21-8 CMF C2 H4 O



=> d 157 abs ibib hitstr 2

ANSWER 2 OF 3 CAPLUS COPYRIGHT 2005 ACS on STN The title amines, giving cured compns. forming clear films, are prepared AB from compds. bearing ≥2 NH2 groups, alkylene carbonates, aliphatic carbonyl compds., and compds. bearing ≥2 OH-reactive groups in specified ratios. Adding 1.0 mol ethylene carbonate over 1 h to 1.0 mol 1,6-hexanediamine in PhMe at 40-50°, heating at 50° until the amine number was <158, adding 2.0 mol MEK and 0.5 g MeC6H4SO3H, distilling H2O, adding 0.5 mol polypropylene glycol-based diepoxy resin (epoxy content 3125 mmol/kg), and heating at 120° until the epoxy content was <50 mmol/kg gave a blocked amine. Use of 76 g this product with 50 g aqueous epoxy resin (Beckopox EP 116) composition (epoxy content 5076 mmol/kg) in the formulation of a clear coating film is exemplified.

1999:736300 CAPLUS ACCESSION NUMBER:

DOCUMENT NUMBER: 131:337884

TITLE: Blocked amines as curing agent for aqueous

one-component epoxy resin compositions

Feola, Roland; Muller, Friedrich; Gmoser, Johann INVENTOR (S): PATENT ASSIGNEE(S): Vianova Resins AG, Austria; Surface Specialties

Austria GM

SOURCE: Eur. Pat. Appl., 11 pp.

CODEN: EPXXDW

DOCUMENT TYPE:

Patent German

LANGUAGE:

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
EP 957121	A2	19991117	EP 1999-108940	19990505
EP 957121	A3	20020522		
EP 957121	B1	20040204		
R: AT, BE, CH,	DE, DK	, ES, FR, GE	B, GR, IT, LI, LU, NL,	SE, MC, PT,
IE, SI, LT,	LV, FI	, RO		
AT 9800802	A	20001015	AT 1998-802	19980512
AT 407748	В	20010525		

AT 258944 E 20040215 AT 1999-108940 19990505 US 6207733 B1 20010327 US 1999-307974 19990510 PRIORITY APPLN. INFO.: AT 1998-802 A 19980512 EP 1999-108940 A 19990505

IT 694-83-7D, 1,2-Cyclohexanediamine, reaction products with alkylene carbonates, **ketones** and **epoxy** resins

1761-71-3D, Bis (4-aminocyclohexyl) methane, reaction products with alkylene carbonates, ketones and epoxy resins

RL: MOA (Modifier or additive use); USES (Uses)

(blocked amines as curing agent for aqueous one-component epoxy resin compns.)

RN 694-83-7 CAPLUS

CN 1,2-Cyclohexanediamine (6CI, 7CI, 8CI, 9CI) (CA INDEX NAME)

RN 1761-71-3 CAPLUS

CN Cyclohexanamine, 4,4'-methylenebis- (9CI) (CA INDEX NAME)

=> d 157 abs ibib hitstr 3

L57 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2005 ACS on STN

AB The title liquid materials contain (meth)acryloyl compound and C2-24 (cyclo)aliphatic primary amines [or their adducts (mol. weight 350-1500) with epoxides, isocyanates, or unsatd. carbonyl compds.] blocked by C≤10 aldehydes or ketones. Thus, a polymer solution (from xylene 3640, glycidyl methacrylate 1560, styrene 2807, and Bu acrylate 1872 g) 4252, acrylic acid 320.4, Cr(III) 2-ethylhexanoate 2.7, and hydroquinone 4.57 g were mixed, aerated at 110° to acid value <2, and mixed with 150 g xylene to give a 53.1% solution with acid value 0.9 and viscosity 170 cPa-s. An equiequiv. mixture of this solution and blocked amine (from dipropylenetriamine 196.5, iso-BuCOMe 330, and Epikote-828 297.7 g) was coated to 40 μ (dry basis) on steel to give a coating with tack-free time 2 h, Persoz hardness 104 and 213 after 1 and 7 days, and good resistance to MEK and gasoline.

ACCESSION NUMBER: 1987:215612 CAPLUS

DOCUMENT NUMBER: 106:215612

TITLE: Fast-drying discoloration-free coating materials

PATENT ASSIGNEE(S): AKZO N. V., Neth.

SOURCE: Jpn. Kokai Tokkyo Koho, 11 pp.

CODEN: JKXXAF

DOCUMENT TYPE: Patent
LANGUAGE: Japanese

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO. KIND DATE APPLICATION NO. DATE

JP	61252273			A2	19861110	JP	1986-68829		19860328
JP	05053186			B4	19930809				
EP	203296			A1	19861203	EP	1986-103844		19860320
EP	203296			B1	19890927				
EP	203296			B2	19970903				
	R: AT,	BE,	CH,	DE,	FR, GB, IT,	LI, LU	J, NL, SE		
AT	46712			E	19891015	AT	1986-103844		19860320
DK	8601385			Α	19860930	DK	1986-1385		19860325
DK	172448			В1	19980713				
МО	8601222			Α	19860930	NO	1986-1222		19860325
NO	175864			В	19940912				
NO	175864			C	19941221				
BR	8601354			Α	19861202	BR	1986-1354		19860326
ES	553476			A1	19870601	ES	1986-553476		19860326
CA	1268889			A1	19900508	CA	1986-505213		19860326
CN	86101964			Α	19860924	CN	1986-101964		19860327
CN	1010318			В	19901107				
FI	8601346			Α	19860930	FI	1986-1346		19860327
FI	80468			В	19900228				
FI	80468			C	19900611				,
AU	8655366			A1	19861002	AU	1986-55366		19860327
ΔU	584689			B2	19890601				
ZA	8602359			Α	19861126	ZA	1986-2359		19860327
US	4990577			Α	19910205	US	1989-309449		19890210
PRIORIT	Y APPLN.	INFO	. :			NL	1985-952	Α	19850329
						EP	1986-103844	Α	19860320
						US	1986-844412	B1	19860326

IT 6864-37-5D, 3,3'-Dimethyl-4,4'-diaminodicyclohexylmethane, reaction products with ketones

RL: USES (Uses)

(crosslinking agents for room temperature-curable acrylic coatings)

RN 6864-37-5 CAPLUS

CN Cyclohexanamine, 4,4'-methylenebis[2-methyl- (9CI) (CA INDEX NAME)

$$\begin{array}{c|c} \text{Me} & \\ \text{H}_2\text{N} & \\ \text{Me} & \\ \end{array}$$

=>

---Logging off of STN---

=>

Executing the logoff script...

=> LOG Y

COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	86.76	2830.87
DIGGOIDE MOIDES (FOR OUR LEVING ACCOUNTS)	SINCE FILE	TOTAL
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	DINCE FIDE	CECCION

CA SUBSCRIBER PRICE

-6.57 -6.57

STN INTERNATIONAL LOGOFF AT 16:28:37 ON 15 FEB 2005